

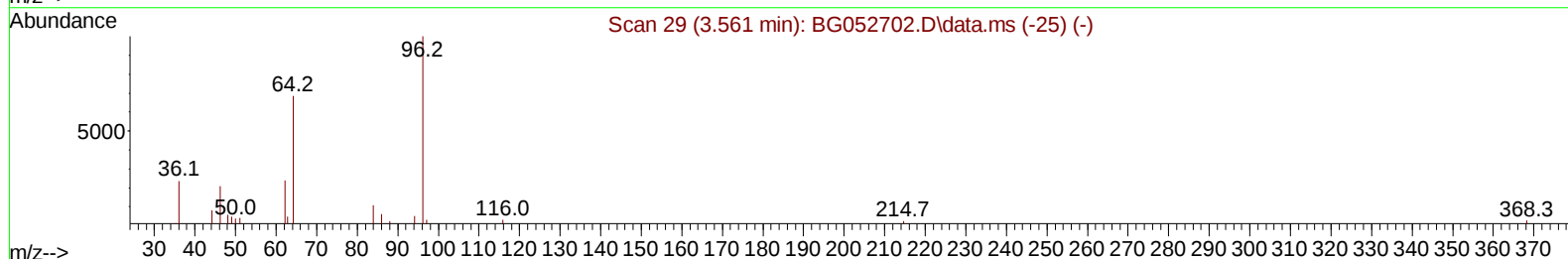
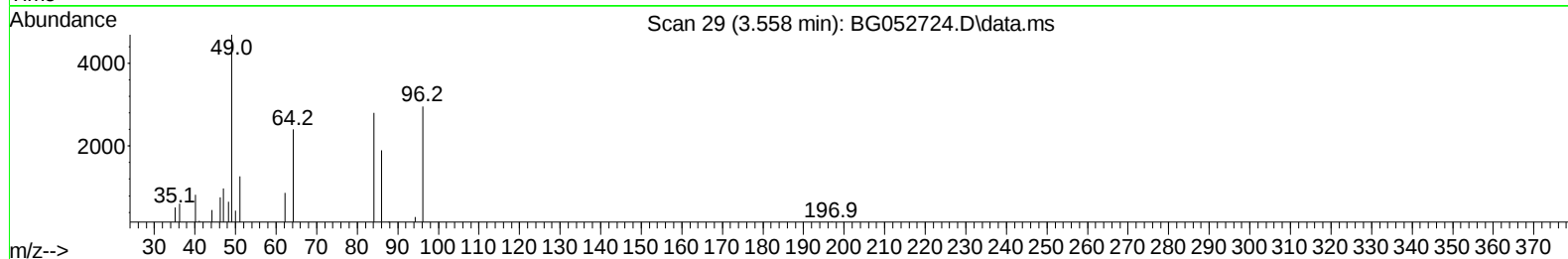
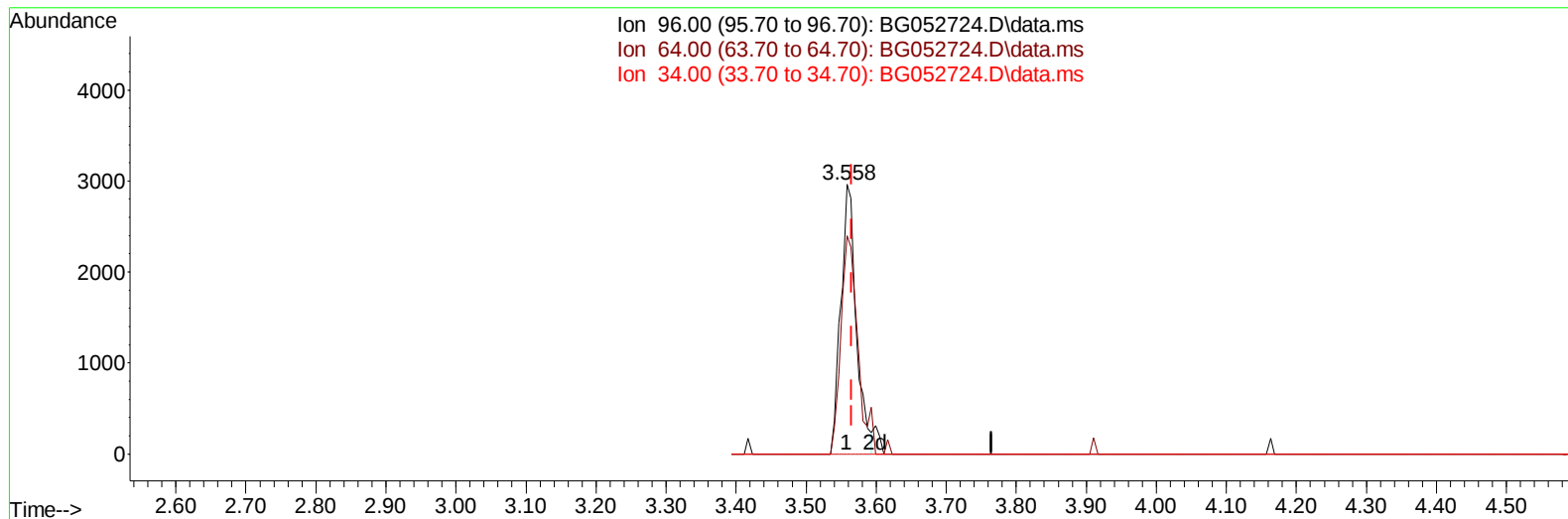
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031822\
Data File : BG052724.D
Acq On : 19 Mar 2022 00:12
Operator : CG/JU
Sample : N1882-14
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBRY4

Manual IntegrationsAPPROVED

Quant Time: Mar 19 01:01:28 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 18 00:55:42 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/21/2022
Supervised By :Yogesh Patel 03/24/2022



TIC: BG052724.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.558min (-0.007) 5.88 ng/uL

response 4602

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	80.99
34.00	0.00	0.00
0.00	0.00	0.00

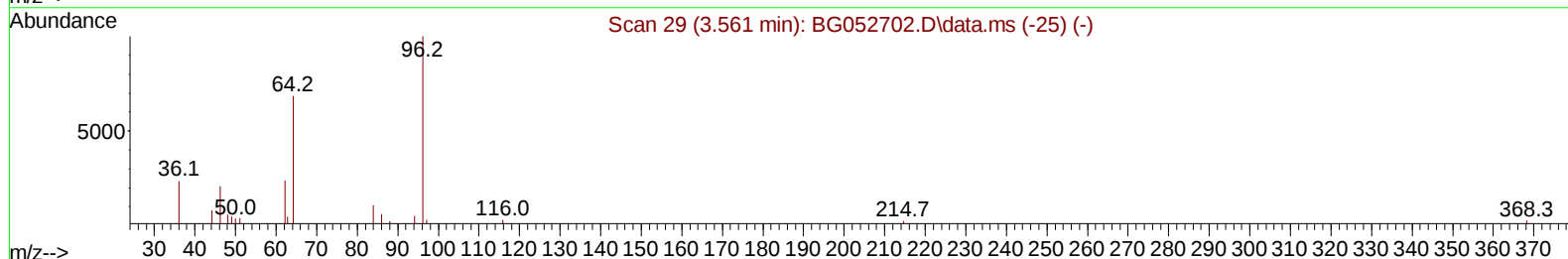
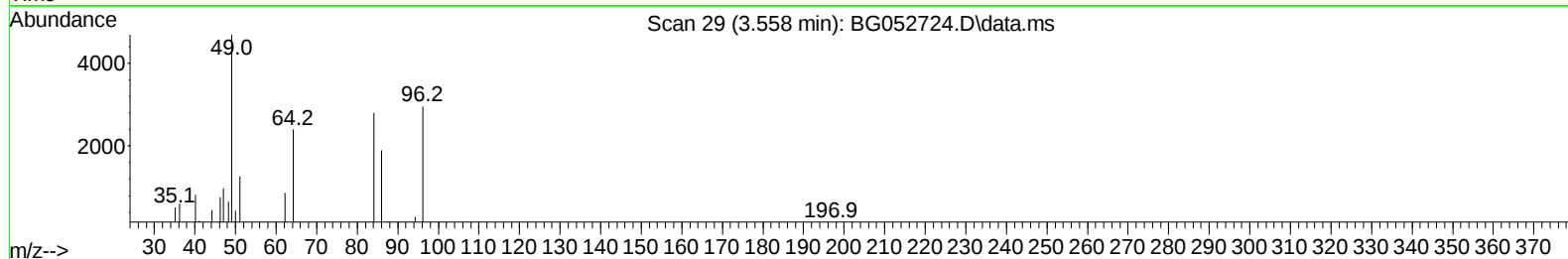
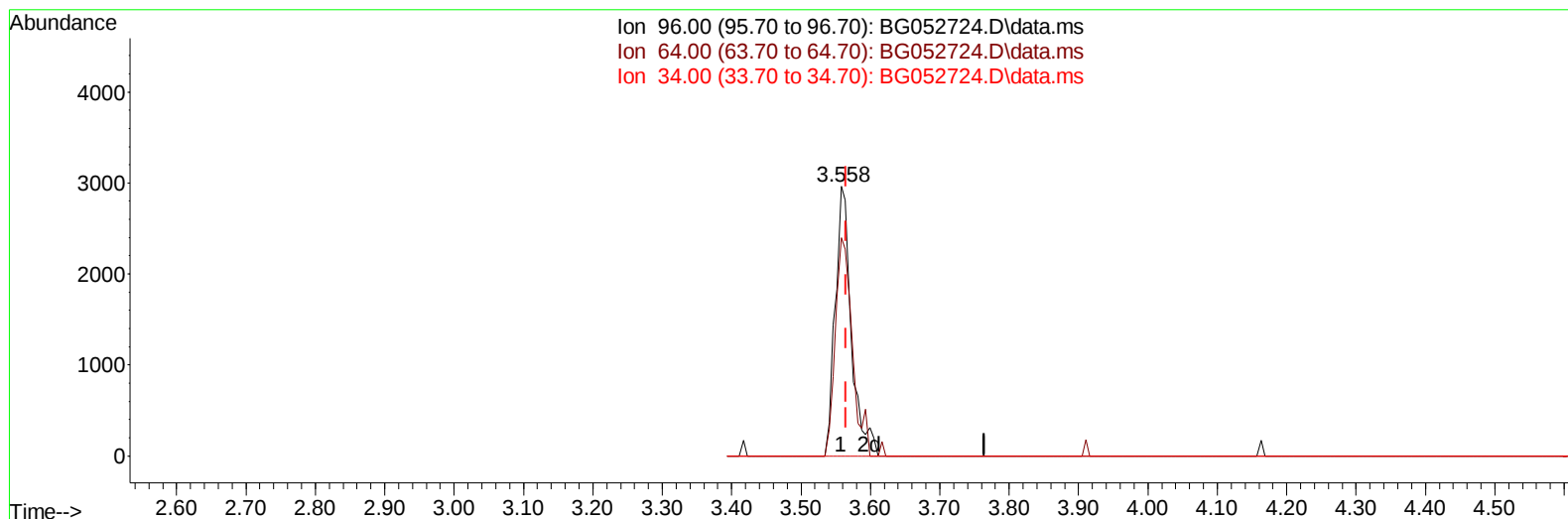
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TIC: BG052724.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.558min (-0.007) 6.11 ng/uL m

response 4780

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	80.99
34.00	0.00	0.00
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.146	152	27641	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.960	136	113334	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.766	164	84641	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.510	188	209488	20.000	ng/ul	0.00
79) Chrysene-d12	21.798	240	221611	20.000	ng/ul	0.00
88) Perylene-d12	25.117	264	214563	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.558	96	4780m	6.108	ng/uL	0.00
4) Pyridine-d5	3.975	84	57215	29.287	ng/ul	0.00
7) Phenol-d5	7.288	99	88451	35.728	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.464	67	55125	36.458	ng/ul	0.00
11) 2-Chlorophenol-d4	7.676	132	60585	36.041	ng/ul	0.00
15) 4-Methylphenol-d8	8.839	113	65488	36.517	ng/ul	0.00
21) Nitrobenzene-d5	9.303	128	30979	37.973	ng/ul	0.00
24) 2-Nitrophenol-d4	10.032	143	33660	39.117	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.572	165	65940	35.225	ng/ul	0.00
31) 4-Chloroaniline-d4	11.083	131	92283	37.299	ng/ul	0.00
46) Dimethylphthalate-d6	14.167	166	247294	38.031	ng/ul	0.00
49) Acenaphthylene-d8	14.461	160	292178	36.978	ng/ul	0.00
54) 4-Nitrophenol-d4	14.931	143	37620	34.410	ng/ul	0.00
60) Fluorene-d10	15.759	176	216445	37.910	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.859	200	33271	31.209	ng/ul	0.00
73) Anthracene-d10	17.610	188	358376	36.738	ng/ul	0.00
81) Pyrene-d10	19.889	212	450125	36.308	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.888	264	414500	37.388	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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